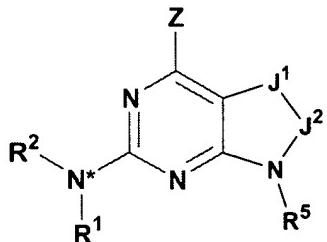


AMENDMENTS TO THE CLAIMS

This listing of claims will replace all prior versions and listings of claims in the application.

Listing of claims:

1. (Currently Amended) A compound of Formula (I)

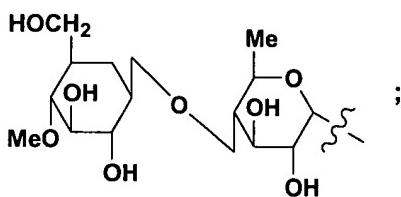


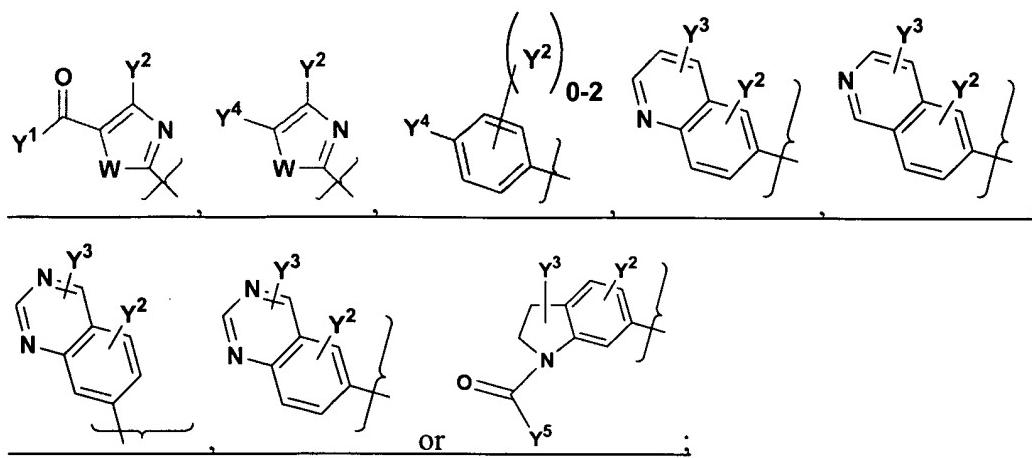
their enantiomers, diastereomers, and pharmaceutically acceptable salts, prodrugs and solvates thereof, wherein:

R¹ is hydrogen or alkyl;

R² is

- (a) heteroaryl or heterocyclo, either of which may be optionally independently substituted with one to three groups selected from T¹, T² and/or T³;
 - (b) aryl substituted with one to three groups selected from T¹, T², and/or T³ provided that at least one of T¹, T² and/or T³ is other than H; or
 - (c) aryl fused to a heteroaryl or heterocyclo ring forming a fused ring system bound to N* through the aryl wherein the fused ring system may be optionally independently substituted with one to three groups selected from T¹, T² and/or T³;
- provided that R² is not





W is O or S;

Y¹ is -NHT¹⁵ or OT¹⁰;

Y² and Y³ are independently hydrogen, halo, OT¹⁰, haloalkyl, or alkyl;

Y⁴ is optionally substituted heteroaryl, cyano, C(O)_tT¹⁰ or S(O)_tNT¹⁴T¹⁵;

Y⁵ is alkyl, NHT¹⁵ or OT¹⁰;

Z is -NR³R⁴, -NR³SO₂R⁶, OR⁴, SR⁴, haloalkyl or halogen;

J¹ is O or S;

J² is optionally substituted C₂alkylene;

R³ and R⁴ are independently H, alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl, cycloalkyl, (cycloalkyl)alkyl, heterocyclo or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁴, T⁵ and/or T⁶;

or R³ and R⁴ may be taken together with the nitrogen atom to which they are attached to form a heterocyclo or heteroaryl ring, either of which is optionally independently substituted where valance allows with one to three groups independently selected from T⁴, T⁵ and/or T⁶;

R⁵ is

- (i) H, cyano, alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or
- (ii) -C(O)_tR⁷, -C(O)-C(O)-C(O)OR⁷ or -SO₂R⁸;

R⁶ is alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl, cycloalkyl, (cycloalkyl)alkyl, heterocylo, or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁴, T⁵ and/or T⁶;

R⁷ is

- (i) H, alkyl, alkenyl, heterocyclo, (heterocyclo)alkyl, (hydroxy)alkyl, (alkoxy)alkyl, (aryloxy)alkyl, heteroaryl, aryl or (aryl)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or
- (ii) $\text{T}^{11}-\text{NHS(O)(T}^{11})-\text{NR}^9\text{R}^{10}$ or $(\text{NR}^9\text{R}^{10})\text{alkyl}$;

R⁸ is

- (i) alkyl, alkenyl, heterocyclo, (heterocyclo)alkyl, (hydroxy)alkyl, (alkoxy)alkyl, (aryloxy)alkyl, heteroaryl, aryl or (aryl)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or
- (ii) $-\text{NR}^9\text{R}^{10}$ or $(\text{NR}^9\text{R}^{10})\text{alkyl}$;

R⁹ and R¹⁰ are independently H, alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl, cycloalkyl, (cycloalkyl)alkyl, heterocylo or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹;

T¹-T⁹ are each independently

- (i)-(iii) alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, $-\text{OT}^{10}$, $-\text{SH}$, $-\text{ST}^{10}$, $-\text{C(O)}_t\text{H}$, $-\text{C(O)}_t\text{T}^{10}$, $-\text{O-C(O)T}^{10}$, $-\text{T}^{17}\text{C(O)}_t\text{N(T}^{11})\text{T}^{10}$, $-\text{SO}_3\text{H}$, $-\text{S(O)}_t\text{T}^{10}$, $-\text{S(O)}_t\text{N(T}^{11})\text{T}^{10}$, $-\text{T}^{12}-\text{NT}^{14}\text{T}^{15}$, $-\text{T}^{12}-\text{N(T}^{11})\text{-T}^{13}-\text{NT}^{14}\text{T}^{15}$, $-\text{T}^{12}-\text{N(T}^{16})\text{-T}^{13}-\text{T}^{10}$, and $-\text{T}^{12}-\text{N(T}^{16})\text{-T}^{13}-\text{H}$; or

(ii) (iv) halo, cyano, nitro, OH, oxo, -SH, amino, $-OT^{10}$, $-ST^{10}$, $-C(O)_tH$, $-C(O)_tT^{10}$, $-O-C(O)T^{10}$, $-T^{17}C(O)_tN(T^{11})T^{10}-$, $-SO_3H$, $-S(O)_tT^{10}$, $-S(O)_tN(T^{11})T^{10}-$, $-T^{12}-NT^{14}T^{15}$, $-T^{12}-N(T^{11})-T^{13}-NT^{14}T^{15}$, $-T^{12}-N(T^{16})-T^{13}-T^{10}-$; or $-T^{12}-N(T^{16})-T^{13}-H$;

t is 1 or 2;

T^{10} is alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl;

T^{12} and T^{13} are each independently a single bond, $-T^{17}-S(O)_t-T^{18}-$, $-T^{17}-C(O)-T^{18}-$, $-T^{17}-C(S)-T^{18}-$, $-T^{17}-O-T^{18}-$, $-T^{17}-S-T^{18}-$, $-T^{17}-O-C(O)-T^{18}-$, $-T^{17}-C(O)_tT^{18}-$, $-T^{17}-C(=NT^{19})-T^{18}-$ or $-T^{17}-C(O)-C(O)-T^{18}-$;

T^{11} , T^{14} , T^{15} , T^{16} and T^{19} are each independently

- (i) hydrogen, alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted where valence permits by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl, (heteroaryl)alkyl, $-SH$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$ and $-S(O)_tT^{22}$; or
- (ii) halo, cyano, nitro, OH, oxo, $-SH$, amino, $-OT^{22}$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$, $-SO_3H$, or $-S(O)_tT^{22}$; or
- (iii) T^{14} and T^{15} may together be alkylene or alkenylene, completing a 3- to 8-membered saturated or unsaturated ring together with the atoms to which they are attached, which ring is substituted with one or more groups listed in the description of T^{20} ; or
- (iv) T^{14} or T^{15} , together with T^{11} , may be alkylene or alkenylene completing a 3- to 8-membered saturated or unsaturated ring together with the nitrogen atoms to which they are attached, which ring is substituted with one or more groups listed in the description of T^{20} ; or
- (v) T^{14} and T^{15} or T^{11} and T^{16} together with the nitrogen atom to which they are attached may combine to form a group $-N=CT^{20}T^{21}$;

T^{17} and T^{18} are each independently a single bond, alkylene, alkenylene or alkynylene; T^{20} and T^{21} are each

(i)[i.] independently hydrogen, alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted where valence permits by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl, (heteroaryl)alkyl, —SH, —ST²², —C(O)_tH, —C(O)_tT²², —O—C(O)T²² and —S(O)_tT²²; or

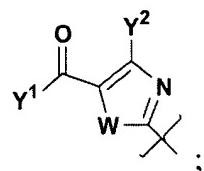
(ii)[ii.] halo, cyano, nitro, OH, oxo, —SH, amino, —OT²², —ST²², —C(O)_tH, —C(O)_tT²², —O—C(O)T²², —SO₃H, or —S(O)_tT²² or —S(O)_tN(T²²)₂; and

T^{22} is alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl.

2. - 3. (Canceled)

4. (Previously Amended) A compound of claim 3, their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein:

R^2 is



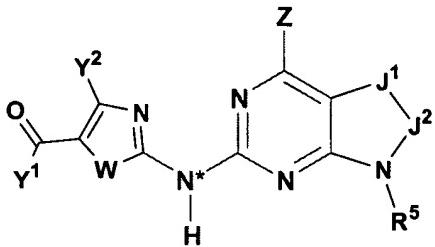
W is O or S;

Y^1 is $-\text{NHT}^{15}$ or OT^{10} ; and

Y^2 is alkyl or haloalkyl.

5. (Canceled)

6. (Currently Amended) A compound of Formula (Ia)



(Ia)

their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein:

W is O or S;

Y¹ is -NHT¹⁵ or OT¹⁰;

Y² is alkyl or haloalkyl;

Z is -NR³R⁴, -NHCH₂CH₂NHC(O)CH₃, or halogen;

J¹ is O;

J² is optionally substituted C₂alkylene;

R³ and R⁴ are independently H, alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl,

cycloalkyl, (cycloalkyl)alkyl, heterocyclo or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁴, T⁵ and/or T⁶;

or R³ and R⁴ may be taken together with the nitrogen atom to which they are attached to form a heterocyclo or heteroaryl ring, either of which is optionally independently substituted where valance allows with one to three groups independently selected from T⁴, T⁵ and/or T⁶;

R⁵ is

(i) H, cyano, alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or

(ii) -C(O)_tR⁷, -C(O)-C(O)-C(O)OR⁷ or -SO₂R⁸;

R⁶ is alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl, cycloalkyl, (cycloalkyl)alkyl, heterocyclo or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁴, T⁵ and/or T⁶;

R⁷ is

- (i) H, alkyl, alkenyl, heterocyclo, (heterocyclo)alkyl, (hydroxy)alkyl, (alkoxy)alkyl, (aryloxy)alkyl, heteroaryl, aryl or (aryl)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or
- (ii) -NR⁹R¹⁰ or (NR⁹R¹⁰)alkyl;

R⁸ is

- (i) alkyl, alkenyl, heterocyclo, (heterocyclo)alkyl, (hydroxy)alkyl, (alkoxy)alkyl, (aryloxy)alkyl, heteroaryl, aryl or (aryl)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹; or
- (ii) -NR⁹R¹⁰ or (NR⁹R¹⁰)alkyl;

R⁹ and R¹⁰ are independently H, alkyl, alkenyl, aryl, (aryl)alkyl, heteroaryl, (heteroaryl)alkyl, cycloalkyl, (cycloalkyl)alkyl, heterocylo or (heterocyclo)alkyl, any of which may be optionally independently substituted where valance allows with one to three groups T⁷, T⁸ and/or T⁹;

T¹-T⁹ are each independently

- (i) alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, -OT¹⁰, -SH, -ST¹⁰, -C(O)_tH, -C(O)_tT¹⁰, -O-C(O)T¹⁰, -T¹⁷C(O)_tN(T¹¹)T¹⁰-, -SO₃H, -S(O)_tT¹⁰, -S(O)_tN(T¹¹)T¹⁰-, -T¹²-NT¹⁴T¹⁵, -T¹²-N(T¹¹)-T¹³-NT¹⁴T¹⁵, -T¹²-N(T¹⁶)-T¹³-T¹⁰, and -T¹²-N(T¹⁶)-T¹³-H;
- (ii) halo, cyano, nitro, OH, oxo, -SH, amino, -OT¹⁰, -ST¹⁰, -C(O)_tH, -C(O)_tT¹⁰, -O-C(O)T¹⁰, -T¹⁷C(O)_tN(T¹¹)T¹⁰-, -SO₃H, -S(O)_tT¹⁰, -S(O)_tN(T¹¹)T¹⁰-, -T¹²-NT¹⁴T¹⁵, -T¹²-N(T¹¹)-T¹³-NT¹⁴T¹⁵, -T¹²-N(T¹⁶)-T¹³-T¹⁰, or -T¹²-N(T¹⁶)-T¹³-H;

t is 1 or 2;

T^{10} is alkyl, **haloalkyl**, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl;

T^{12} and T^{13} are each independently a single bond, $-T^{17}-S(O)_t-T^{18}-$, $-T^{17}-C(O)-T^{18}-$, $-T^{17}-C(S)-T^{18}-$, $-T^{17}-O-T^{18}-$, $-T^{17}-S-T^{18}-$, $-T^{17}-O-C(O)-T^{18}-$, $-T^{17}-C(O)_t-T^{18}-$, $-T^{17}-C(=NT^{19})-T^{18}-$ or $-T^{17}-C(O)-C(O)-T^{18}-$;

T^{11} , T^{14} , T^{15} , T^{16} and T^{19} are each independently

- (i) hydrogen, alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted where valence permits by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl, (heteroaryl)alkyl, $-SH$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$, $-O-C(O)T^{22}$ and $-S(O)T^{22}$; or
- (ii) halo, cyano, nitro, OH, oxo, $-SH$, amino, $-OT^{22}$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$, $-SO_3H$, $-S(O)_tT^{22}$ or $-S(O)_tN(T^{11})T^{22}$; or
- (iii) T^{14} and T^{15} may together be alkylene or alkenylene, completing a 3- to 8-membered saturated or unsaturated ring together with the atoms to which they are attached, which ring is substituted with one or more groups listed in the description of T^{20} ; or
- (iv) T^{14} or T^{15} , together with T^{11} , may be alkylene or alkenylene completing a 3- to 8-membered saturated or unsaturated ring together with the nitrogen atoms to which they are attached, which ring is substituted with one or more groups listed in the description of T^{20} ; or
- (v) T^{14} and T^{15} or T^{11} and T^{16} together with the nitrogen atom to which they are attached may combine to form a group $-N=CT^{20}T^{21}$;

T^{17} and T^{18} are each independently a single bond, alkylene, alkenylene or alkynylene;

T^{20} and T^{21} are each

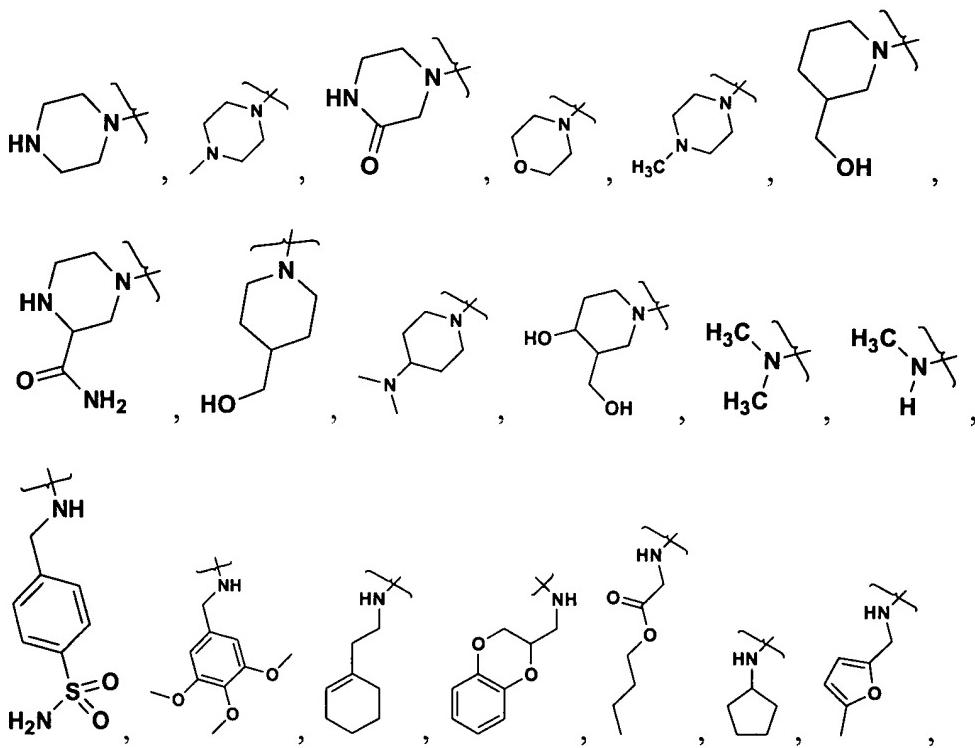
- (i) independently hydrogen, alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl,

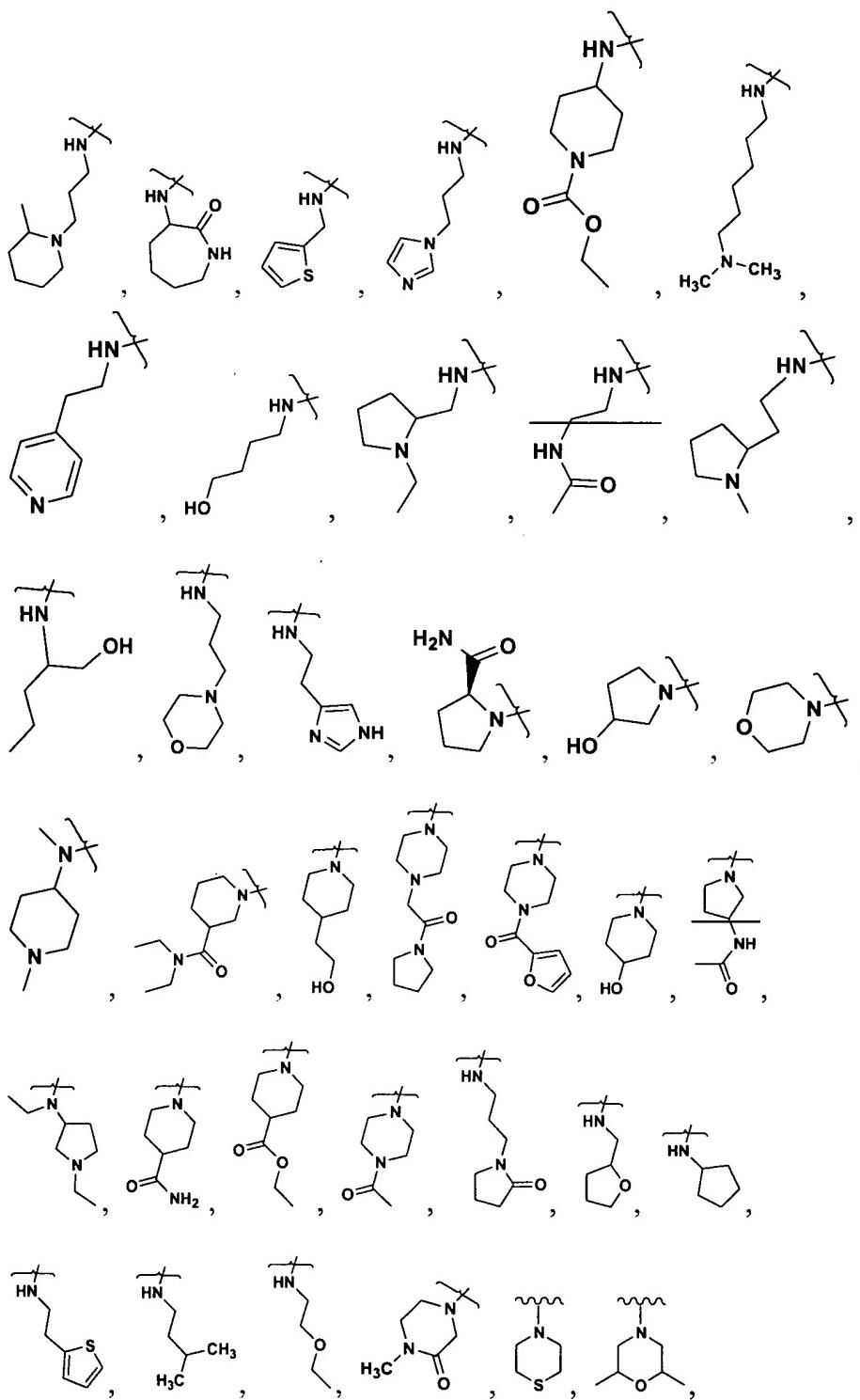
heterocyclo, (heterocylco)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted where valence permits by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyco)alkyl, heteroaryl, (heteroaryl)alkyl, —SH, —ST²², —C(O)_tH, -C(O)_tT²², —O-C(O)T²² and -S(O)_tT²²; or

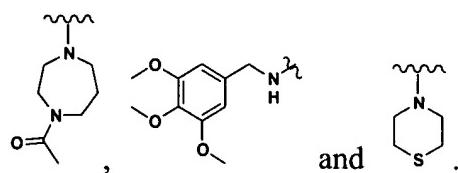
- (ii) halo, cyano, nitro, OH, oxo, -SH, amino, $-OT^{22}$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$, $-SO_3H$, $-S(O)_tT^{22}$ or $S(O)_tN(T^{11})T^{22}$; and

T^{22} is alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl.

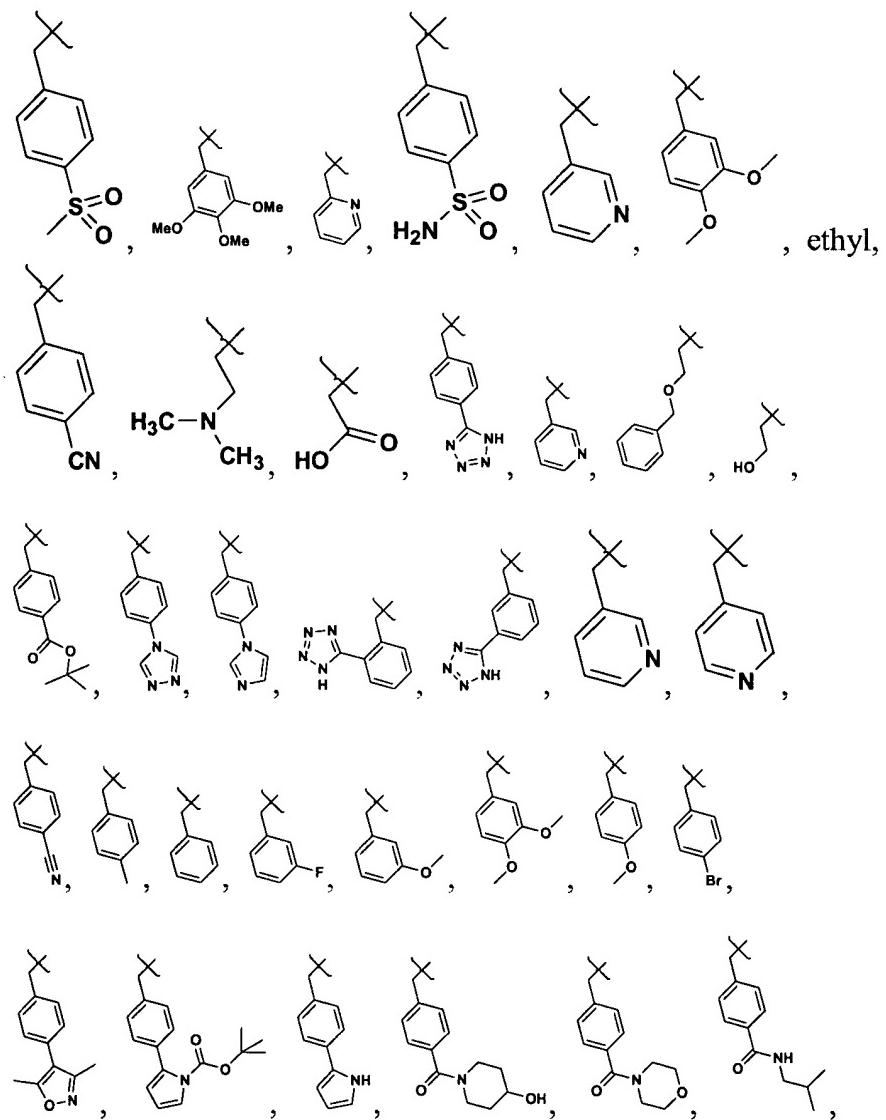
7. (Currently Amended) A compound of claim 6, their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein Z is selected from:

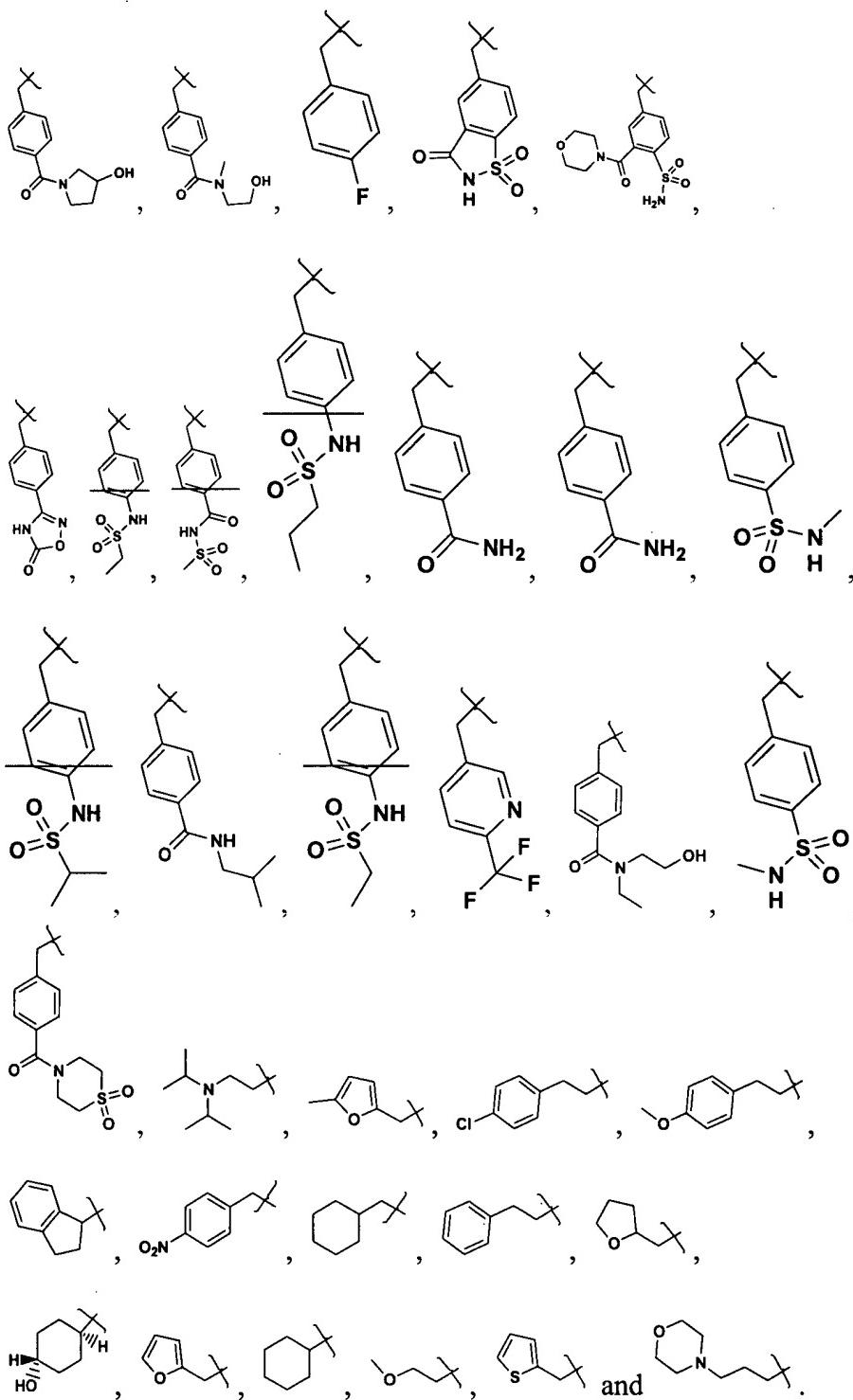




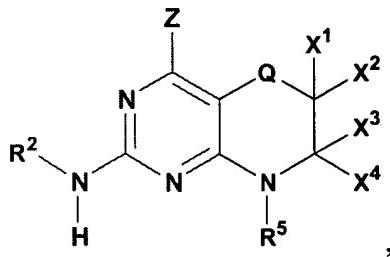


8. (Currently Amended) A compound of claim 6, their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein R⁵ is selected from:





9. (Currently Amended) A compound of claim 1 having Formula (II)



(II)

their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein:

Q is O or S; and

X¹, X², X³ and X⁴ are

- (i) independently chosen from hydrogen, T¹⁰, OT¹⁰ and NT¹⁴T¹⁵; or
- (ii) X¹ and X² or X³ and X⁴ may be taken together to be a carbonyl group.

10. (Currently Amended) A compound of claim 9, their enantiomers, diastereomers, and pharmaceutically acceptable salts, and solvates thereof, wherein Q is O.

11. – 14. (Canceled)

15. (Previously Amended) A pharmaceutical composition comprising at least one compound of claim 1 and a pharmaceutically acceptable carrier or diluent.

16. (Currently Amended) The pharmaceutical composition of claim 15 comprising a pharmaceutically acceptable carrier or diluent and at least one compound selected from:

- i-(i) 2-[8-(4-Methanesulfonyl-benzyl)-4-(3-oxo-piperazin-1-yl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester;
- 4-Methyl-2-[4-morpholin-4-yl-8-(3,4,5-trimethoxy-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester;
- 4-Methyl-2-[4-morpholin-4-yl-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester;

2-[4-(4-Hydroxy-piperidin-1-yl)-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester;
4-Methyl-2-[4-(3-oxo-piperazin-1-yl)-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester;
2-[8-(4-Methanesulfonyl-benzyl)-4-morpholin-4-yl-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester; and
2-[4-(4-Hydroxy-piperidin-1-yl)-8-(4-methanesulfonyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester; or

ii-(ii) the enantiomers, diastereomers, and pharmaceutically acceptable salts, prodrugs and solvates of each of (i).

17. – 19 (Canceled).

20. (Currently Amended) A method of treating a leukocyte activation-associated disorder which comprises administering an effective amount of at least one compound of claim 1, 6, or 21 wherein said disorder is transplant rejection, graft versus host disease, rheumatoid arthritis, multiple sclerosis, juvenile diabetes, asthma, inflammatory bowel disease, ischemic or reperfusion injury, cell proliferation, or psoriasis.

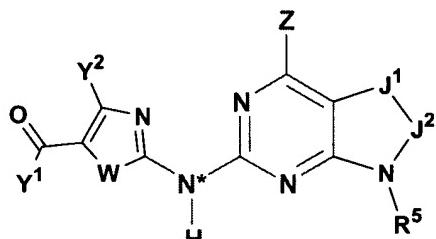
21. (Currently Amended) A compound selected from

i.(i) 2-[8-(4-Methanesulfonyl-benzyl)-4-(3-oxo-piperazin-1-yl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester;
4-Methyl-2-[4-morpholin-4-yl-8-(3,4,5-trimethoxy-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester;
4-Methyl-2-[4-morpholin-4-yl-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester;
2-[4-(4-Hydroxy-piperidin-1-yl)-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester;

4-Methyl-2-[4-(3-oxo-piperazin-1-yl)-8-(4-sulfamoyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-thiazole-5-carboxylic acid ethyl ester; 2-[8-(4-Methanesulfonyl-benzyl)-4-morpholin-4-yl-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester; and 2-[4-(4-Hydroxy-piperidin-1-yl)-8-(4-methanesulfonyl-benzyl)-6,7-dihydro-pyrimido[5,4-b][1,4]oxazin-2-ylamino]-4-methyl-thiazole-5-carboxylic acid ethyl ester; or

ii.(ii) the enantiomers, diastereomers, and pharmaceutically acceptable salts, prodrugs and solvates of each of (i).

22. (Currently Amended) A compound of Formula (Ia)



(Ia)

their enantiomers, diastereomers, pharmaceutically acceptable salts, and solvates thereof, wherein:

W is O or S;

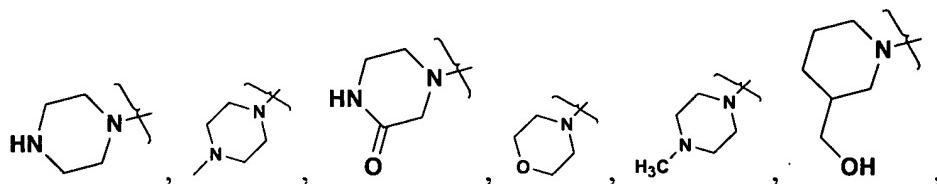
Y¹ is -NHT¹⁵ or OT¹⁰;

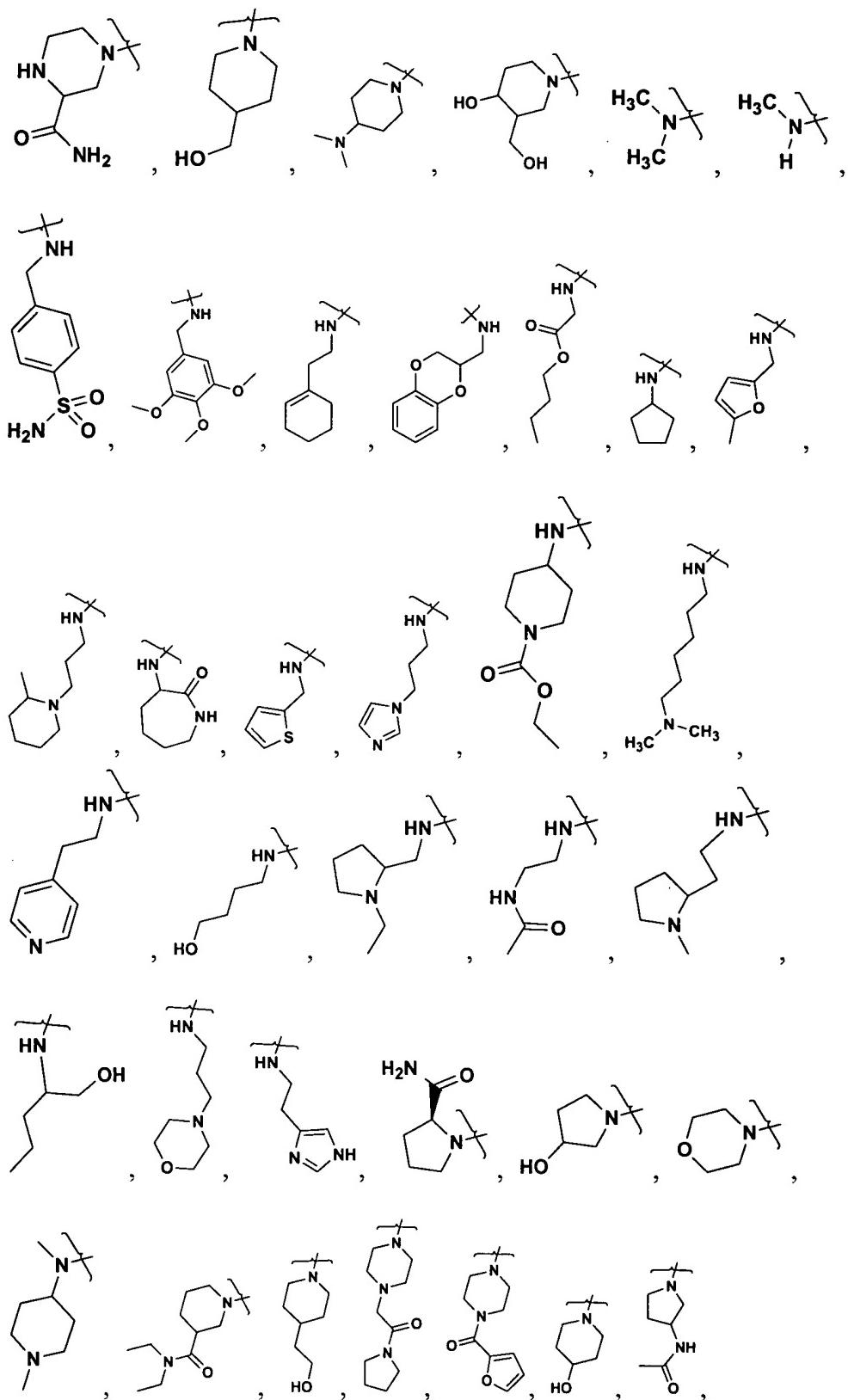
Y² is alkyl or haloalkyl;

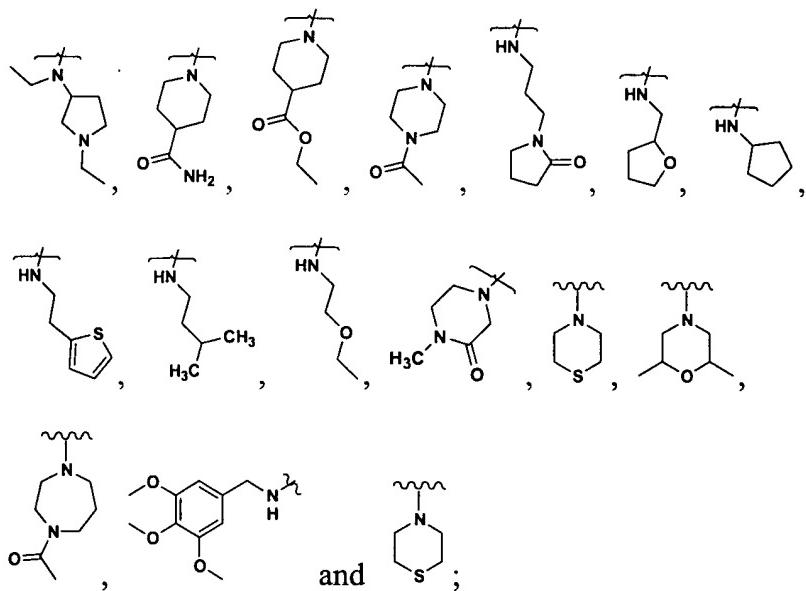
J¹ is O;

J² is optionally substituted C₂alkylene;

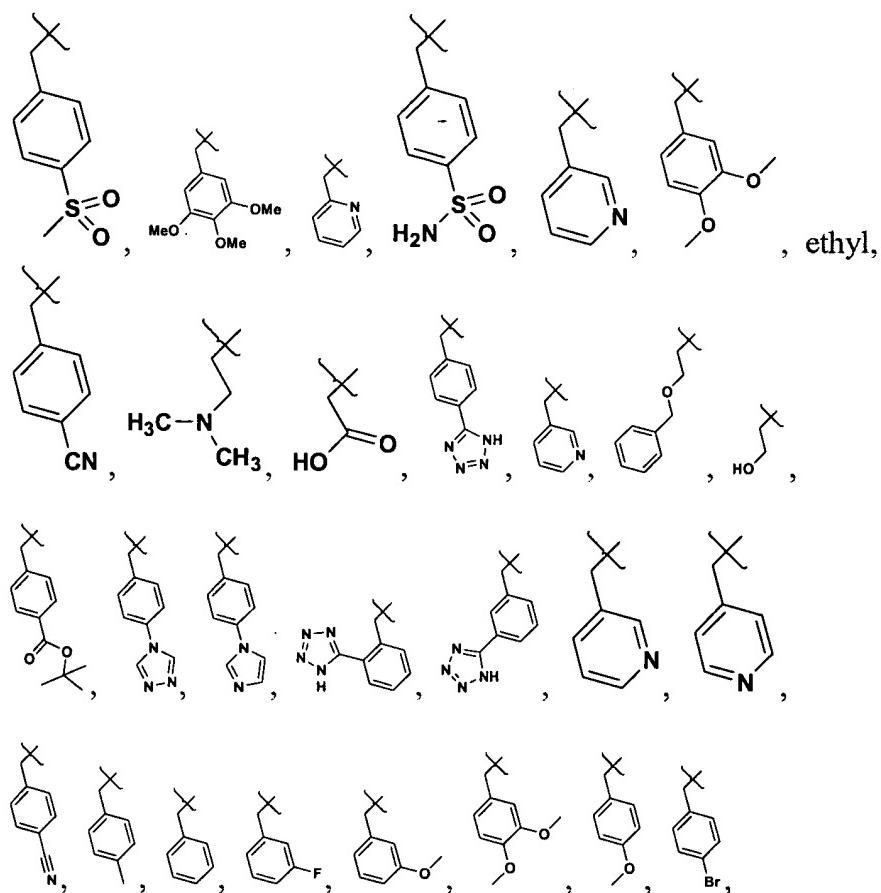
Z is selected from:

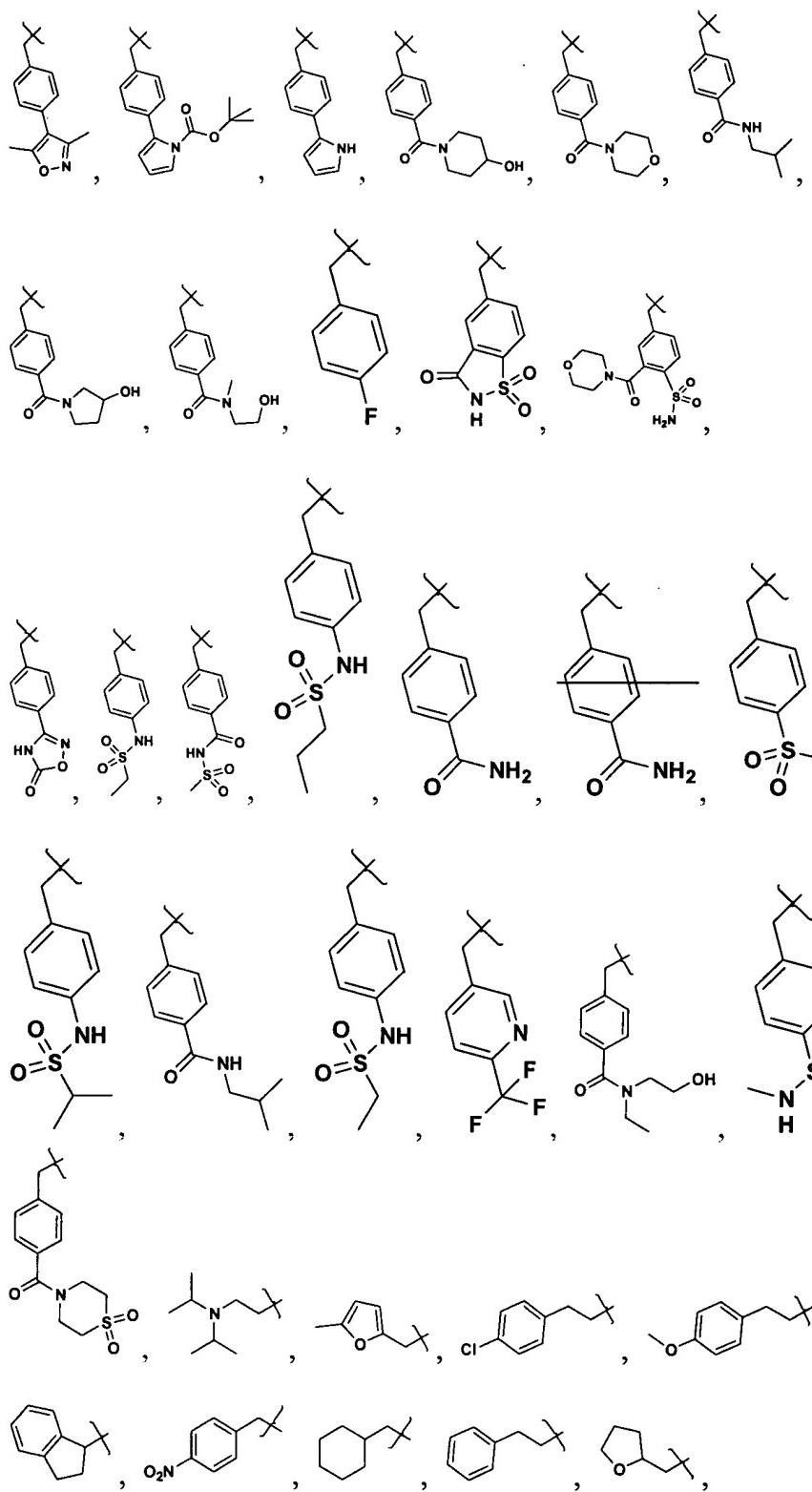


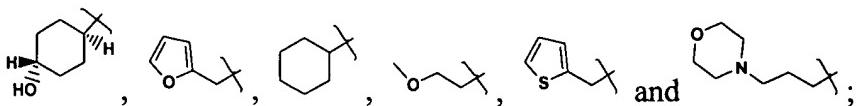




R^5 is selected from:







T^{10} is alkyl, haloalkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl;

T^{15} is (i) hydrogen, alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl, any of which may be optionally independently substituted where valence permits by one or more groups selected from alkyl, (hydroxy)alkyl, halo, cyano, nitro, OH, oxo, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl, (heteroaryl)alkyl, $-SH$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$ and $-S(O)_tT^{22}$; or
(ii) halo, cyano, nitro, OH, oxo, $-SH$, amino, $-OT^{22}$, $-ST^{22}$, $-C(O)_tH$, $-C(O)_tT^{22}$, $-O-C(O)T^{22}$, $-SO_3H$, or $-S(O)_tT^{22}$ or $S(O)_tN(T^{14})T^{22}$;

t is 1 or 2; and

T^{22} is alkyl, (hydroxy)alkyl, (alkoxy)alkyl, alkenyl, alkynyl, cycloalkyl, (cycloalkyl)alkyl, cycloalkenyl, (cycloalkenyl)alkyl, aryl, (aryl)alkyl, heterocyclo, (heterocyclo)alkyl, heteroaryl or (heteroaryl)alkyl.